

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111824\
 Data File : BM048778.D
 Acq On : 18 Nov 2024 16:02
 Operator : RC/JU
 Sample : P4829-10
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 E29W4

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 11/19/2024
 Supervised By :mohammad ahmed 11/20/2024

Quant Time: Nov 18 16:39:30 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM110624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 18 11:56:05 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.629	152	3026m	0.400	ng/u1	-0.12
4) Naphthalene-d8	10.404	136	8951	0.400	ng/u1	-0.08
9) Acenaphthene-d10	14.276	164	4875	0.400	ng/u1	-0.02
13) Phenanthrene-d10	17.024	188	9959m	0.400	ng/u1	-0.05
17) Chrysene-d12	21.243	240	7915	0.400	ng/u1	0.00
23) Perylene-d12	23.423	264	5013m	0.400	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.085	96	14315	3.921	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.032	152	2983	0.258	ng/u1	-0.06
18) Fluoranthene-d10	19.059	212	8726	0.393	ng/u1	-0.03

Target Compounds	Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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